

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** NASA JPL  
**Collection Date:** December 21, 2000  
**LDC Report Date:** March 1, 2001  
**Matrix:** Air  
**Parameters:** Volatile Halogenated/Aromatic Hydrocarbons  
**Validation Level:** EPA Level III  
**Laboratory:** HP Labs  
**Sample Delivery Group (SDG):** 2K1221W1

**Sample Identification**

SWW35-VPA-038  
SWW35-VPD-039  
SWW34-VPA-040  
SWW34-VPB-041  
SWW34-VPB-042DUP  
SWW34-VPD-043  
SWW34-VPF-044  
SWW34-VPG-045  
SWW34-VPH-046  
SWW32-VPB-047  
SWW32-VPB-048DUP  
SWW32-VPC-049

## Introduction

This data review covers 12 air samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Methods 8010 and 8020 for Volatile Halogenated/Aromatic Hydrocarbons.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review (October 1999) as there are no current guidelines for the method stated above.

A table summarizing all data qualification flags is provided at the end of this report. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blank results are summarized in Section III.

Field duplicates are summarized in Section IX.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- J Indicates an estimated value.
- R Quality control indicates the data is not usable.
- N Presumptive evidence of presence of the constituent.
- UU Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration of compounds was performed for the primary (quantitation) column and confirmation column as required by these methods.

The percent relative standard deviations (%RSD) were less than or equal to 20.0% for all compounds.

### **b. Calibration Verification**

Calibration verification was performed at the required frequencies.

The percent differences (%D) of calibration factors in continuing standard mixtures were within the 15.0% QC limits.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No volatile halogenated/aromatic hydrocarbon contaminants were found in the method blanks.

## **IV. Accuracy and Precision Data**

### **a. Surrogate Recovery**

Surrogates were added to all samples and blanks as required by the methods. All surrogate recoveries (%R) were within QC limits.

### **b. Matrix Spike/Matrix Spike Duplicates**

Matrix spike (MS) and matrix spike duplicate (MSD) analyses were not required by the method.

### **c. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits.

## V. Target Compound Identification

Raw data were not reviewed for this SDG.

## VI. Compound Quantitation and CRQLs

Raw data were not reviewed for this SDG.

## VII. System Performance

Raw data were not reviewed for this SDG.

## VIII. Overall Assessment of Data

Data flags are summarized at the end of this report.

## IX. Field Duplicates

Samples SVW34-VPB-041 and SVW34-VPB-042DUP and samples SVW32-VPB-047 and SVW32-VPB-048DUP were identified as field duplicates. No volatile halogenated/aromatic hydrocarbons were detected in any of the samples with the following exceptions:

Compound	Concentration (ug/L)		RPD
	SVW34-VPB-041	SVW34-VPB-042DUP	
Carbon tetrachloride	5.9	5.9	0

## X. Field Blanks

No field blanks were identified in this SDG.

**NASA JPL**

**Volatile Halogenated/Aromatic Hydrocarbons - Data Qualification Summary - SDG 2K1221W1**

No Sample Data Qualified in this SDG

**NASA JPL**

**Volatile Halogenated/Aromatic Hydrocarbons - Laboratory Blank Data Qualification Summary - SDG 2K1221W1**

No Sample Data Qualified in this SDG

GEOFON PROJECT # 04-4304-480

JPL

4800 OAK GROVE DRIVE

PASADENA, CA

HP Labs Project #2K1221W1

GC SHIMADZU 14A FRONT

VOLATILE HALOGENATED AND AROMATIC HYDROCARBONS (EPA Method 8010/8020) ANALYSES OF SOIL VAPOR

SOIL VAPOR DATA IN UG/L-VAPOR

	BLANK	SVW35-VPA-038	SVW35-VPD-039	SVW34-VPA-040	SVW34-VPB-041	SVW34-VPB-042 DUP	SVW34-VPD-043
DATE	12/21/00	12/21/00	12/21/00	12/21/00	12/21/00	12/21/00	12/21/00
SAMPLING TIME	05:40	06:59	07:25	08:17	08:37	09:05	09:30
ANALYSIS TIME	05:42	07:04	07:28	08:19	08:46	09:10	09:33
SAMPLING DEPTH (feet)	--	20	60	20	35	35	65
VOLUME WITHDRAWN (cc)	200	80	240	80	140	140	260
VOLUME INJECTED	1	1	1	1	1	1	1
DILUTION FACTOR	1	1	1	1	1	1	1
CARBON TETRACHLORIDE	nd	nd	nd	nd	5.9	5.9	nd
CHLOROETHANE/BROMOMETHANE	nd	nd	nd	nd	nd	nd	nd
CHLOROFORM	nd	nd	nd	nd	nd	nd	nd
1,1-DICHLORO ETHANE	nd	nd	nd	nd	nd	nd	nd
1,2-DICHLORO ETHANE	nd	nd	nd	nd	nd	nd	nd
1,1-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd	nd
CIS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd	nd
TRANS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd	nd
DICHLOROMETHANE	nd	nd	nd	nd	nd	nd	nd
TETRACHLORO ETHENE	nd	nd	nd	nd	nd	nd	nd
1,1,1,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd	nd	nd
1,1,2,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd	nd	nd
1,1,1-TRICHLORO ETHANE	nd	nd	nd	nd	nd	nd	nd
1,1,2-TRICHLORO ETHANE	nd	nd	nd	nd	nd	nd	nd
TRICHLORO ETHENE	nd	nd	nd	nd	nd	nd	nd
VINYL CHLORIDE	nd	nd	nd	nd	nd	nd	nd
TRICHLOROFLUOROMETHANE (FR11)	nd	nd	nd	nd	nd	nd	nd
DICHLORODIFLUOROMETHANE (FR12)	nd	nd	nd	nd	nd	nd	nd
1,1,2-TRICHLOROTRIFLUOROETHANE (FR113)	nd	nd	nd	nd	nd	nd	nd
BENZENE	nd	nd	nd	nd	nd	nd	nd
ETHYLBENZENE	nd	nd	nd	nd	nd	nd	nd
TOLUENE	nd	nd	nd	nd	nd	nd	nd
m&p-XYLENES	nd	nd	nd	nd	nd	nd	nd
o-XYLENE	nd	nd	nd	nd	nd	nd	nd
SURROGATES							
1,4 DIFLUORO BENZENE	107%	105%	101%	105%	106%	107%	97%
CHLOROBENZENE	97%	104%	100%	108%	106%	108%	96%
4 BROMOFLUORO BENZENE	99%	106%	103%	109%	108%	112%	100%

ND INDICATES NOT DETECTED AT A DETECTION LIMIT OF 1.0 UG/L-VAPOR FOR EACH COMPOUND

ANALYSES PERFORMED ON-SITE IN DOHS CERTIFIED MOBILE LABORATORY (CERT #1745)

ANALYSES PERFORMED BY: ALLEN GLOVER

DATA REVIEWED BY: JAMES E. PICKER

3/2/01

GEOFON PROJECT # 04-4304-480

JPL

4800 OAK GROVE DRIVE

PASADENA, CA

HP Labs Project #2K1221W1

GC SHIMADZU 14A FRONT

VOLATILE HALOGENATED AND AROMATIC HYDROCARBONS (EPA Method 8010/8020) ANALYSES OF SOIL VAPOR

SOIL VAPOR DATA IN UG/L-VAPOR

	SVW34-VPF-044	SVW34-VPG-045	SVW34-VPH-046	SVW32-VPB-047	SVW32-VPB-048 DUP	SVW32-VPC-049
DATE	12/21/00	12/21/00	12/21/00	12/21/00	12/21/00	12/21/00
SAMPLING TIME	09:55	10:15	10:43	11:07	11:30	11:55
ANALYSIS TIME	09:57	10:21	10:45	11:10	11:34	11:58
SAMPLING DEPTH (feet)	95	108	118	40	40	55
VOLUME WITHDRAWN (cc)	380	435	475	160	160	220
VOLUME INJECTED	1	1	1	1	1	1
DILUTION FACTOR	1	1	1	1	1	1
CARBON TETRACHLORIDE	nd	nd	26	nd	nd	nd
CHLOROETHANE/BROMOMETHANE	nd	nd	nd	nd	nd	nd
CHLOROFORM	nd	nd	2.4	nd	nd	nd
1,1-DICHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,2-DICHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,1-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd
CIS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd
TRANS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd
DICHLOROMETHANE	nd	nd	nd	nd	nd	nd
TETRACHLORO ETHENE	nd	nd	nd	nd	nd	nd
1,1,1,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,1,2,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,1,1-TRICHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,1,2-TRICHLORO ETHANE	nd	nd	nd	nd	nd	nd
TRICHLORO ETHENE	nd	nd	nd	nd	nd	nd
VINYL CHLORIDE	nd	nd	nd	nd	nd	nd
TRICHLOROFLUOROMETHANE (FR11)	nd	nd	nd	nd	nd	nd
DICHLORODIFLUOROMETHANE (FR12)	nd	nd	nd	nd	nd	nd
1,1,2-TRICHLOROTRIFLUOROETHANE (FR113)	nd	nd	1.4	nd	nd	nd
BENZENE	nd	nd	nd	nd	nd	nd
ETHYLBENZENE	nd	nd	nd	nd	nd	nd
TOLUENE	nd	nd	nd	nd	nd	nd
m&p-XYLENES	nd	nd	nd	nd	nd	nd
o-XYLENE	nd	nd	nd	nd	nd	nd
SURROGATES						
1,4 DIFLUORO BENZENE	95%	95%	98%	102%	94%	93%
CHLOROBENZENE	97%	97%	100%	100%	95%	95%
4 BROMOFLUORO BENZENE	100%	99%	100%	104%	97%	96%

ND INDICATES NOT DETECTED AT A DETECTION LIMIT OF 1.0 UG/L-VAPOR FOR EACH COMPOUND

ANALYSES PERFORMED ON-SITE IN DOHS CERTIFIED MOBILE LABORATORY (CERT #1745)

ANALYSES PERFORMED BY: ALLEN GLOVER

DATA REVIEWED BY: JAMES E. PICKER

3/2/01

LDC #: 6038D23 **VALIDATION COMPLETENESS WORKSHEET**SDG #: 2K1221W1 X EPA Level III    NFESC Level C

Laboratory: HP Labs

Date: 2/28/01

Page: 1 of 1

Reviewer: ET

2nd Reviewer: N

**METHOD:** GC Volatile Halogenated/Aromatic Hydrocarbons (EPA SW 846 Method 8010/8020)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 12/21/00
IIa.	Initial calibration	A	% RSD
IIb.	Calibration verification	A	% D
III.	Blanks	A	
IVa.	Surrogate recovery	A	
IVb.	Matrix spike/Matrix spike duplicates	N	
IVc.	Laboratory control samples	PT/A	
V.	Target compound identification	N	
VI.	Compound Quantitation and CRQLs	N	
VII.	System Performance	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	SW	D = 4 + 5 * D <sub>1</sub> = 10 + 11 * = 20
X.	Field blanks	N	

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:

Air

1	SVW35-VPA-038	11	SVW32-VPB-048DUP	21		31	
2	SVW35-VPD-039	12	SVW32-VPC-049	22		32	
3	SVW34-VPA-040	13	BLK	23		33	
4	SVW34-VPB-041	14		24		34	
5	SVW34-VPB-042DUP	15		25		35	
6	SVW34-VPD-043	16		26		36	
7	SVW34-VPF-044	17		27		37	
8	SVW34-VPG-045	18		28		38	
9	SVW34-VPH-046	19		29		39	
10	SVW32-VPB-047	20		30		40	

Notes:



LDC #: 6038023  
SDG #: 2K/22/W/

# TARGET COMPOUND WORKSHEET

Page: 1 of 1  
Reviewer: FT  
2nd Reviewer: me

METHOD: VOA (EPA SW 846 Method 8240/8260/8021))

A. Chloromethane*	P. Bromodichloromethane	EE. Ethylbenzene**	TT. 1,2-Dibromoethane	III. n-Butylbenzene
B. Bromomethane	Q. 1,2-Dichloropropane**	FF. Styrene	UU. 1,1,1,2-Tetrachloroethane	JJJ. 1,2-Dichlorobenzene
C. Vinyl chloride**	R. cis-1,3-Dichloropropene	GG. Xylene, total	VV. Isopropylbenzene	KKK. 1,2,4-Trichlorobenzene
D. Chloroethane	S. Trichloroethene	HH. Vinyl acetate	WW. Bromobenzene	LLL. Hexachlorobutadiene
E. Methylene chloride	T. Dibromochloromethane	II. 2-Chloroethylvinyl ether	XX. 1,2,3-Trichloropropane	MMM. Naphthalene
F. Acetone	U. 1,1,2-Trichloroethane	JJ. Dichlorodifluoromethane	YY. n-Propylbenzene	NNN. 1,2,3-Trichlorobenzene
G. Carbon disulfide	V. Benzene	KK. Trichlorofluoromethane	ZZ. 2-Chlorotoluene	OOO. 1,3,5-Trichlorobenzene
H. 1,1-Dichloroethene**	W. trans-1,3-Dichloropropene	LL. Methyl-tert-butyl ether	AAA. 1,3,5-Trimethylbenzene	PPP. trans-1,2-Dichloroethene
I. 1,1-Dichloroethane*	X. Bromoform*	MM. 1,2-Dibromo-3-chloropropane	BBB. 4-Chlorotoluene	QQQ. cis-1,2-Dichloroethene
J. 1,2-Dichloroethene	Y. 4-Methyl-2-pentanone	NN. Diethyl ether	CCC. tert-Butylbenzene	RRR.
K. Chloroform**	Z. 2-Hexanone	OO. 2,2-Dichloropropane	DDD. 1,2,4-Trimethylbenzene	SSS.
L. 1,2-Dichloroethane	AA. Tetrachloroethene	PP. Bromochloromethane	EEE. sec-Butylbenzene	TTT.
M. 2-Butanone	BB. 1,1,2,2-Tetrachloroethane*	QQ. 1,1-Dichloropropene	FFF. 1,3-Dichlorobenzene	UUU.
N. 1,1,1-Trichloroethane	CC. Toluene**	RR. Dibromomethane	GGG. p-Isopropyltoluene	VVV.
O. Carbon tetrachloride	DD. Chlorobenzene*	SS. 1,3-Dichloropropane	HHH. 1,4-Dichlorobenzene	WWW.

\* = System performance check compounds (SPCC) for RF ; \*\* = Calibration check compounds (CCC) for %RSD.

Notes: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

LDC #: 6038 D23  
SDG #: 2K/22/W/

**VALIDATION FINDINGS WORKSHEET**  
**Field Duplicates**

Page: 1 of 1  
Reviewer: FT  
2nd reviewer:

**METHOD:** GC Volatiles (EPA SW 846 Method 8010/8020)

☒ Y ☐ N ☐ N/A  
☒ Y ☐ N ☐ N/A

Were field duplicate pairs identified in this SDG?

Were target compounds detected in the field duplicate pairs?

Compound	Concentration ( <u>ug/L</u> )		RPD
	4	5	
0	5.9	5.9	0

Compound	Concentration ( )		RPD

Compound	Concentration ( )		RPD

Compound	Concentration ( )		RPD

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** NASA JPL  
**Collection Date:** December 22, 2000  
**LDC Report Date:** March 1, 2001  
**Matrix:** Air  
**Parameters:** Volatile Halogenated/Aromatic Hydrocarbons  
**Validation Level:** EPA Level III  
**Laboratory:** HP Labs

**Sample Delivery Group (SDG):** 2K1222W1

**Sample Identification**

SVW32-VPD-050  
SVW32-VPE-051  
SVW32-VPH-052  
SVW32-VPI-053  
SVW32-VPI-054DUP  
SVW32-VPJ-055  
SVW39-VPA-056  
SVW39-VPC-057  
SVW39-VPD-058  
SVW39-VPE-059  
SVW39-VPE-060DUP  
SVW39-VPF-061

## Introduction

This data review covers 12 air samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Methods 8010 and 8020 for Volatile Halogenated/Aromatic Hydrocarbons.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review (October 1999) as there are no current guidelines for the method stated above.

A table summarizing all data qualification flags is provided at the end of this report. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blank results are summarized in Section III.

Field duplicates are summarized in Section IX.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- J Indicates an estimated value.
- R Quality control indicates the data is not usable.
- N Presumptive evidence of presence of the constituent.
- UU Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration of compounds was performed for the primary (quantitation) column and confirmation column as required by these methods.

The percent relative standard deviations (%RSD) were less than or equal to 20.0% for all compounds.

### **b. Calibration Verification**

Calibration verification was performed at the required frequencies.

The percent differences (%D) of calibration factors in continuing standard mixtures were within the 15.0% QC limits.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No volatile halogenated/aromatic hydrocarbon contaminants were found in the method blanks.

## **IV. Accuracy and Precision Data**

### **a. Surrogate Recovery**

Surrogates were added to all samples and blanks as required by the methods. All surrogate recoveries (%R) were within QC limits.

### **b. Matrix Spike/Matrix Spike Duplicates**

Matrix spike (MS) and matrix spike duplicate (MSD) analyses were not required by the method.

### **c. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits.

## V. Target Compound Identification

Raw data were not reviewed for this SDG.

## VI. Compound Quantitation and CRQLs

Raw data were not reviewed for this SDG.

## VII. System Performance

Raw data were not reviewed for this SDG.

## VIII. Overall Assessment of Data

Data flags are summarized at the end of this report.

## IX. Field Duplicates

Samples SVW32-VPI-053 and SVW32-VPI-054DUP and samples SVW39-VPE-059 and SVW39-VPE-060DUP were identified as field duplicates. No volatile halogenated/aromatic hydrocarbons were detected in any of the samples with the following exceptions:

Compound	Concentration (ug/L)		RPD
	SVW39-VPE-059	SVW39-VPE-060DUP	
Carbon disulfide	2.4	2.1	13
Trichloroethene	1.3	1.3	0
1,1,2-Trichlorotrifluoroethane	17	16	6

## X. Field Blanks

No field blanks were identified in this SDG.

**NASA JPL**

**Volatile Halogenated/Aromatic Hydrocarbons - Data Qualification Summary - SDG 2K1222W1**

No Sample Data Qualified in this SDG

**NASA JPL**

**Volatile Halogenated/Aromatic Hydrocarbons - Laboratory Blank Data Qualification Summary - SDG 2K1222W1**

No Sample Data Qualified in this SDG

GEOFON PROJECT # 04-4304-480

JPL

4800 OAK GROVE DRIVE

PASADENA, CA

HP Labs Project #2K1222W1

GC SHIMADZU 14A FRONT

VOLATILE HALOGENATED AND AROMATIC HYDROCARBONS (EPA Method 8010/8020) ANALYSES OF SOIL VAPOR

SOIL VAPOR DATA IN UG/L-VAPOR

	BLANK	SVW32-VPD-050	SVW32-VPE-051	SVW32-VPH-052	SVW32-VPI-053	SVW32-VPI-054 DUP	SVW32-VPJ-055
DATE	12/22/00	12/22/00	12/22/00	12/22/00	12/22/00	12/22/00	12/22/00
SAMPLING TIME	05:53	06:54	07:19	07:44	08:08	08:31	08:55
ANALYSIS TIME	05:54	06:58	07:22	07:45	08:09	08:33	08:58
SAMPLING DEPTH (feet)	--	70	90	155	180	180	195
VOLUME WITHDRAWN (cc)	200	280	360	620	720	720	780
VOLUME INJECTED	1	1	1	1	1	1	1
DILUTION FACTOR	1	1	1	1	1	1	1
CARBON TETRACHLORIDE	nd	nd	nd	14	nd	nd	nd
CHLOROETHANE/BROMOMETHANE	nd	nd	nd	nd	nd	nd	nd
CHLOROFORM	nd	nd	nd	nd	nd	nd	nd
1,1-DICHLORO ETHANE	nd	nd	nd	nd	nd	nd	nd
1,2-DICHLORO ETHANE	nd	nd	nd	nd	nd	nd	nd
1,1-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd	nd
CIS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd	nd
TRANS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd	nd
DICHLOROMETHANE	nd	nd	nd	nd	nd	nd	nd
TETRACHLORO ETHENE	nd	nd	nd	nd	nd	nd	nd
1,1,1,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd	nd	nd
1,1,2,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd	nd	nd
1,1,1-TRICHLORO ETHANE	nd	nd	nd	nd	nd	nd	nd
1,1,2-TRICHLORO ETHANE	nd	nd	nd	nd	nd	nd	nd
TRICHLORO ETHENE	nd	nd	nd	nd	nd	nd	nd
VINYL CHLORIDE	nd	nd	nd	nd	nd	nd	nd
TRICHLOROFLUOROMETHANE (FR11)	nd	nd	nd	nd	nd	nd	nd
DICHLORODIFLUOROMETHANE (FR12)	nd	nd	nd	nd	nd	nd	nd
1,1,2-TRICHLOROTRIFLUOROETHANE (FR113)	nd	nd	nd	17	nd	nd	nd
BENZENE	nd	nd	nd	nd	nd	nd	nd
ETHYLBENZENE	nd	nd	nd	nd	nd	nd	nd
TOLUENE	nd	nd	nd	nd	nd	nd	nd
m&p-XYLENES	nd	nd	nd	nd	nd	nd	nd
o-XYLENE	nd	nd	nd	nd	nd	nd	nd
SURROGATES							
1,4 DIFLUORO BENZENE	105%	96%	87%	93%	94%	97%	102%
CHLOROBENZENE	95%	98%	88%	93%	93%	97%	105%
4 BROMOFLUORO BENZENE	97%	102%	88%	95%	95%	100%	107%

ND INDICATES NOT DETECTED AT A DETECTION LIMIT OF 1.0 UG/L-VAPOR FOR EACH COMPOUND

ANALYSES PERFORMED ON-SITE IN DOHS CERTIFIED MOBILE LABORATORY (CERT #1745)

ANALYSES PERFORMED BY: ALLEN GLOVER

DATA REVIEWED BY: JAMES E. PICKER

12/31/01





GEOFON PROJECT # 04-4304-480

JPL

4800 OAK GROVE DRIVE

PASADENA, CA

HP Labs Project #201222W1

GC SHIMADZU 14A FRONT

VOLATILE HALOGENATED AND AROMATIC HYDROCARBONS (EPA Method 8010/8020) ANALYSES OF SOIL VAPOR

SOIL VAPOR DATA IN UG/L-VAPOR

	SVW39-VPA-056	SVW39-VPC-057	SVW39-VPD-058	SVW39-VPE-059	SVW39-VPE-060 DUP	SVW-VPF-061
DATE	12/22/00	12/22/00	12/22/00	12/22/00	12/22/00	12/22/00
SAMPLING TIME	09:20	09:42	10:05	10:30	10:54	11:17
ANALYSIS TIME	09:21	09:45	10:09	10:32	10:56	11:20
SAMPLING DEPTH (feet)	20	50	70	85	85	100
VOLUME WITHDRAWN (cc)	80	200	280	340	340	400
VOLUME INJECTED	1	1	1	1	1	1
DILUTION FACTOR	1	1	1	1	1	1
CARBON TETRACHLORIDE	nd	nd	nd	2.4	2.1	5.0
CHLOROETHANE/BROMOMETHANE	nd	nd	nd	nd	nd	nd
CHLOROFORM	nd	nd	nd	nd	nd	nd
1,1-DICHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,2-DICHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,1-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd
CIS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd
TRANS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd
DICHLOROMETHANE	nd	nd	nd	nd	nd	nd
TETRACHLORO ETHENE	nd	nd	nd	nd	nd	nd
1,1,1,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,1,2,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,1,1-TRICHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,1,2-TRICHLORO ETHANE	nd	nd	nd	nd	nd	nd
TRICHLORO ETHENE	nd	nd	nd	1.3	1.3	2.3
VINYL CHLORIDE	nd	nd	nd	nd	nd	nd
TRICHLOROFLUOROMETHANE (FR11)	nd	nd	nd	nd	nd	nd
DICHLORODIFLUOROMETHANE (FR12)	nd	nd	nd	nd	nd	nd
1,1,2-TRICHLOROTRIFLUOROETHANE (FR113)	nd	nd	nd	17	16	21
BENZENE	nd	nd	nd	nd	nd	nd
ETHYLBENZENE	nd	nd	nd	nd	nd	nd
TOLUENE	nd	nd	nd	nd	nd	nd
m&p-XYLENES	nd	nd	nd	nd	nd	nd
o-XYLENE	nd	nd	nd	nd	nd	nd
SURROGATES						
1,4 DIFLUORO BENZENE	107%	95%	98%	99%	97%	99%
CHLOROBENZENE	109%	97%	98%	102%	94%	98%
4 BROMOFLUORO BENZENE	112%	101%	100%	103%	96%	102%

ND INDICATES NOT DETECTED AT A DETECTION LIMIT OF 1.0 UG/L-VAPOR FOR EACH COMPOUND

ANALYSES PERFORMED ON-SITE IN DOHS CERTIFIED MOBILE LABORATORY (CERT #1745)

ANALYSES PERFORMED BY: ALLEN GLOVER

DATA REVIEWED BY: JAMES E. PICKER

12/31/01

LDC #: 6038E23 **VALIDATION COMPLETENESS WORKSHEET**  
 SDG #: 2K1222W1 X EPA Level III    NFESC Level C  
 Laboratory: HP Labs

Date: 2/28/01  
 Page: 1 of 1  
 Reviewer: FA  
 2nd Reviewer:   

**METHOD:** GC Volatile Halogenated/Aromatic Hydrocarbons (EPA SW 846 Method 8010/8020)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 12/22/00
IIa.	Initial calibration	A	% RSD
IIb.	Calibration verification	A	% D
III.	Blanks	A	
IVa.	Surrogate recovery	A	
IVb.	Matrix spike/Matrix spike duplicates	N	
IVc.	Laboratory control samples	NA <sup>FA</sup>	
V.	Target compound identification	N	
VI.	Compound Quantitation and CRQLs	N	
VII.	System Performance	N	
VIII.	Overall assessment of data	A	
IX.	Field duplicates	SW	*D = 4 & 5      D <sub>1</sub> = 10 & 11      * = ND
X.	Field blanks	N	

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet

ND = No compounds detected  
 R = Rinsate  
 FB = Field blank

D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples:

AIR

1	SVW32-VPD-050	†1	SVW39-VPE-060DUP	D <sub>1</sub>	21		31	
2	SVW32-VPE-051	†2	SVW39-VPF-061		22		32	
3	SVW32-VPH-052	13	BLK		23		33	
4	SVW32-VPI-053	D			24		34	
5	SVW32-VPI-054DUP	D			25		35	
6	SVW32-VPJ-055				26		36	
7	SVW39-VPA-056				27		37	
8	SVW39-VPD-057				28		38	
9	SVW39-VPD-058				29		39	
10	SVW39-VPE-059	D <sub>1</sub>			30		40	

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

LDC #: 6038 E 23  
SDG #: 2F1222 W 1

# TARGET COMPOUND WORKSHEET

Page: 1 of 1  
Reviewer: #7  
2nd Reviewer: n

METHOD: VOA (EPA SW 846 Method 8240/8260/8021))

A. Chloromethane*	P. Bromodichloromethane	EE. Ethylbenzene**	TT. 1,2-Dibromoethane	III. n-Butylbenzene
B. Bromomethane	Q. 1,2-Dichloropropane**	FF. Styrene	UU. 1,1,1,2-Tetrachloroethane	JJJ. 1,2-Dichlorobenzene
C. Vinyl chloride**	R. cis-1,3-Dichloropropene	GG. Xylene, total	VV. Isopropylbenzene	KKK. 1,2,4-Trichlorobenzene
D. Chloroethane	S. Trichloroethene	HH. Vinyl acetate	WW. Bromobenzene	LLL. Hexachlorobutadiene
E. Methylene chloride	T. Dibromochloromethane	II. 2-Chloroethylvinyl ether	XX. 1,2,3-Trichloropropane	MMM. Naphthalene
F. Acetone	U. 1,1,2-Trichloroethane	JJ. Dichlorodifluoromethane	YY. n-Propylbenzene	NNN. 1,2,3-Trichlorobenzene
G. Carbon disulfide	V. Benzene	KK. Trichlorofluoromethane	ZZ. 2-Chlorotoluene	OOO. 1,3,5-Trichlorobenzene
H. 1,1-Dichloroethene**	W. trans-1,3-Dichloropropene	LL. Methyl-tert-butyl ether	AAA. 1,3,5-Trimethylbenzene	PPP. trans-1,2-Dichloroethene
I. 1,1-Dichloroethane*	X. Bromoform*	MM. 1,2-Dibromo-3-chloropropane	BBB. 4-Chlorotoluene	QQQ. cis-1,2-Dichloroethene
J. 1,2-Dichloroethene	Y. 4-Methyl-2-pentanone	NN. Diethyl ether	CCC. tert-Butylbenzene	RRR.
K. Chloroform**	Z. 2-Hexanone	OO. 2,2-Dichloropropane	DDD. 1,2,4-Trimethylbenzene	SSS.
L. 1,2-Dichloroethane	AA. Tetrachloroethene	PP. Bromochloromethane	EEE. sec-Butylbenzene	TTT.
M. 2-Butanone	BB. 1,1,2,2-Tetrachloroethane*	QQ. 1,1-Dichloropropene	FFF. 1,3-Dichlorobenzene	UUU.
N. 1,1,1-Trichloroethane	CC. Toluene**	RR. Dibromomethane	GGG. p-Isopropyltoluene	VVV.
O. Carbon tetrachloride	DD. Chlorobenzene*	SS. 1,3-Dichloropropane	HHH. 1,4-Dichlorobenzene	WWW.

\* = System performance check compounds (SPCC) for RF ; \*\* = Calibration check compounds (CCC) for %RSD.

ZZZ- 1,1,2-Trichlorotrifluoroethane (FR113)

Notes: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

LDC #: 6038 E23  
SDG #: 2K1222 W1

VALIDATION FINDINGS WORKSHEET  
Field Duplicates

Page: 1 of 1  
Reviewer: FT  
2nd reviewer: FT

METHOD: GC Volatiles (EPA SW 846 Method 8010/8020)

Y N N/A  
Y N N/A

Were field duplicate pairs identified in this SDG?

Were target compounds detected in the field duplicate pairs?

Compound	Concentration (ug/L)		RPD
	10	11	
G	2.4	2.1	13
S	1.3	1.3	0
222	17	16	FTX 6

Compound	Concentration ( )		RPD

Compound	Concentration ( )		RPD

Compound	Concentration ( )		RPD

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** NASA JPL  
**Collection Date:** December 27, 2000  
**LDC Report Date:** March 1, 2001  
**Matrix:** Air  
**Parameters:** Volatile Halogenated/Aromatic Hydrocarbons  
**Validation Level:** EPA Level III  
**Laboratory:** HP Labs  
**Sample Delivery Group (SDG):** 2K1227W1

**Sample Identification**

SVW39-VPI-062  
SVW38-VPA-063  
SVW38-VPB-064  
SVW38-VPC-065  
SVW38-VPC-066DUP  
SVW38-VPD-067  
SVW38-VPF-068  
SVW38-VPG-069  
SVW38-VPJ-070

## Introduction

This data review covers 9 air samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Methods 8010 and 8020 for Volatile Halogenated/Aromatic Hydrocarbons.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review (October 1999) as there are no current guidelines for the method stated above.

A table summarizing all data qualification flags is provided at the end of this report. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blank results are summarized in Section III.

Field duplicates are summarized in Section IX.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- J Indicates an estimated value.
- R Quality control indicates the data is not usable.
- N Presumptive evidence of presence of the constituent.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration of compounds was performed for the primary (quantitation) column and confirmation column as required by these methods.

The percent relative standard deviations (%RSD) were less than or equal to 20.0% for all compounds.

### **b. Calibration Verification**

Calibration verification was performed at the required frequencies.

The percent differences (%D) of calibration factors in continuing standard mixtures were within the 15.0% QC limits.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No volatile halogenated/aromatic hydrocarbon contaminants were found in the method blanks.

## **IV. Accuracy and Precision Data**

### **a. Surrogate Recovery**

Surrogates were added to all samples and blanks as required by the methods. All surrogate recoveries (%R) were within QC limits.

### **b. Matrix Spike/Matrix Spike Duplicates**

Matrix spike (MS) and matrix spike duplicate (MSD) analyses were not required by the method.

### **c. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits.

## **V. Target Compound Identification**

Raw data were not reviewed for this SDG.

## **VI. Compound Quantitation and CRQLs**

Raw data were not reviewed for this SDG.

## **VII. System Performance**

Raw data were not reviewed for this SDG.

## **VIII. Overall Assessment of Data**

Data flags are summarized at the end of this report.

## **IX. Field Duplicates**

Samples SVW38-VPC-065 and SVW38-VPC-066DUP were identified as field duplicates. No volatile halogenated/aromatic hydrocarbons were detected in any of the samples.

## **X. Field Blanks**

No field blanks were identified in this SDG.



**NASA JPL**

**Volatile Halogenated/Aromatic Hydrocarbons - Data Qualification Summary - SDG 2K1227W1**

No Sample Data Qualified in this SDG

**NASA JPL**

**Volatile Halogenated/Aromatic Hydrocarbons - Laboratory Blank Data Qualification Summary - SDG 2K1227W1**

No Sample Data Qualified in this SDG

GEOFON PROJECT # 04-4304-480

JPL

4800 OAK GROVE DRIVE

PASADENA, CA

HP Labs Project #2K1227W1

GC SHIMADZU 14A FRONT

VOLATILE HALOGENATED AND AROMATIC HYDROCARBONS (EPA Method 8010/8020) ANALYSES OF SOIL VAPOR

SOIL VAPOR DATA IN UG/L-VAPOR

	BLANK	SVW39-VPI-062	SVW38-VPA-063	SVW36-VPB-064	SVW38-VPC-065
DATE	12/27/00	12/27/00	12/27/00	12/27/00	12/27/00
SAMPLING TIME	06:05	07:00	07:22	07:45	08:10
ANALYSIS TIME	06:05	07:03	07:26	07:50	08:13
SAMPLING DEPTH (feet)	--	130	25	45	65
VOLUME WITHDRAWN (cc)	200	520	100	180	260
VOLUME INJECTED	1	1	1	1	1
DILUTION FACTOR	1	1	1	1	1
CARBON TETRACHLORIDE	nd	2.4	nd	nd	nd
CHLOROETHANE/BROMOMETHANE	nd	nd	nd	nd	nd
CHLOROFORM	nd	nd	nd	nd	nd
1,1-DICHLORO ETHANE	nd	nd	nd	nd	nd
1,2-DICHLORO ETHANE	nd	nd	nd	nd	nd
1,1-DICHLORO ETHENE	nd	nd	nd	nd	nd
CIS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd
TRANS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd
DICHLOROMETHANE	nd	nd	nd	nd	nd
TETRACHLORO ETHENE	nd	nd	nd	nd	nd
1,1,1,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd
1,1,2,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd
1,1,1-TRICHLORO ETHANE	nd	nd	nd	nd	nd
1,1,2-TRICHLORO ETHANE	nd	nd	nd	nd	nd
TRICHLORO ETHENE	nd	5.2	nd	nd	nd
VINYL CHLORIDE	nd	nd	nd	nd	nd
TRICHLOROFLUOROMETHANE (FR11)	nd	nd	nd	nd	nd
DICHLORODIFLUOROMETHANE (FR12)	nd	nd	nd	nd	nd
1,1,2-TRICHLOROTRIFLUOROETHANE (FR113)	nd	2.1	nd	nd	nd
BENZENE	nd	nd	nd	nd	nd
ETHYLBENZENE	nd	nd	nd	nd	nd
TOLUENE	nd	nd	nd	nd	nd
m&p-XYLENES	nd	nd	nd	nd	nd
o-XYLENE	nd	nd	nd	nd	nd
SURROGATES					
1,4 DIFLUORO BENZENE	96%	97%	94%	94%	96%
CHLOROBENZENE	90%	97%	95%	95%	95%
4 BROMOFLUORO BENZENE	91%	100%	98%	97%	98%

ND INDICATES NOT DETECTED AT A DETECTION LIMIT OF 1.0 UG/L-VAPOR FOR EACH COMPOUND

ANALYSES PERFORMED ON-SITE IN DOHS CERTIFIED MOBILE LABORATORY (CERT #1745)

ANALYSES PERFORMED BY: ALLEN GLOVER

DATA REVIEWED BY: JAMES E. PICKER

3/2/01



GEOFON PROJECT # 04-4304-480

JPL

4800 OAK GROVE DRIVE

PASADENA, CA

HP Labs Project #2K1227W1

GC SHIMADZU 14A FRONT

VOLATILE HALOGENATED AND AROMATIC HYDROCARBONS (EPA Method 8010/8020) ANALYSES OF SOIL VAPOR

SOIL VAPOR DATA IN UG/L-VAPOR

	SVW38-VPC-066 DUP	SVW38-VPD-067	SVW38-VPF-068	SVW38-VPG-069	SVW38-VPJ-070
DATE	12/27/00	12/27/00	12/27/00	12/27/00	12/27/00
SAMPLING TIME	08:35	08:57	09:20	09:45	10:14
ANALYSIS TIME	08:37	09:00	09:24	09:48	10:15
SAMPLING DEPTH (feet)	65	80	110	125	170
VOLUME WITHDRAWN (cc)	260	320	440	500	680
VOLUME INJECTED	1	1	1	1	1
DILUTION FACTOR	1	1	1	1	1
CARBON TETRACHLORIDE	nd	nd	3.0	2.3	4.6
CHLOROETHANE/BROMOMETHANE	nd	nd	nd	nd	nd
CHLOROFORM	nd	nd	nd	nd	nd
1,1-DICHLORO ETHANE	nd	nd	nd	nd	nd
1,2-DICHLORO ETHANE	nd	nd	nd	nd	nd
1,1-DICHLORO ETHENE	nd	nd	nd	nd	nd
CIS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd
TRANS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd
DICHLOROMETHANE	nd	nd	nd	nd	nd
TETRACHLORO ETHENE	nd	nd	nd	nd	nd
1,1,1,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd
1,1,2,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd
1,1,1-TRICHLORO ETHANE	nd	nd	nd	nd	nd
1,1,2-TRICHLORO ETHANE	nd	nd	nd	nd	nd
TRICHLORO ETHENE	nd	nd	1.2	1.0	2.2
VINYL CHLORIDE	nd	nd	nd	nd	nd
TRICHLOROFLUOROMETHANE (FR11)	nd	nd	nd	nd	nd
DICHLORODIFLUOROMETHANE (FR12)	nd	nd	nd	nd	nd
1,1,2-TRICHLOROTRIFLUOROETHANE (FR113)	nd	nd	1.5	1.4	5.9
BENZENE	nd	nd	nd	nd	nd
ETHYLBENZENE	nd	nd	nd	nd	nd
TOLUENE	nd	nd	nd	nd	nd
m&p-XYLENES	nd	nd	nd	nd	nd
o-XYLENE	nd	nd	nd	nd	nd
SURROGATES					
1,4 DIFLUORO BENZENE	98%	99%	95%	99%	104%
CHLOROBENZENE	98%	99%	96%	101%	103%
4 BROMOFLUORO BENZENE	100%	102%	98%	102%	107%

ND INDICATES NOT DETECTED AT A DETECTION LIMIT OF 1.0 UG/L-VAPOR FOR EACH COMPOUND

ANALYSES PERFORMED ON-SITE IN DOHS CERTIFIED MOBILE LABORATORY (CERT #1745)

ANALYSES PERFORMED BY: ALLEN GLOVER

DATA REVIEWED BY: JAMES E. PICKER

per 3/2/01

LDC #: 6038F23 **VALIDATION COMPLETENESS WORKSHEET**

SDG #: 2K1227W1 X EPA Level III    NFESC Level C

Laboratory: HP Labs

Date: 2/28/01

Page: 1 of 1

Reviewer: ET

2nd Reviewer: N

**METHOD:** GC Volatile Halogenated/Aromatic Hydrocarbons (EPA SW 846 Method 8010/8020)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 12/27/00
IIa.	Initial calibration	A	% RSD
IIb.	Calibration verification	A	% D
III.	Blanks	A	
IVa.	Surrogate recovery	A	
IVb.	Matrix spike/Matrix spike duplicates	N	
IVc.	Laboratory control samples	FLWA	
V.	Target compound identification	N	
VI.	Compound Quantitation and CRQLs	N	
VII.	System Performance	N	
VIII.	Overall assessment of data		
IX.	Field duplicates	ND	D = 4 & 5
X.	Field blanks	N	

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples:

AIR

1	SVW39-VPI-062	11		21		31	
2	SVW38-VPA-063	12		22		32	
3	SVW38-VPB-064	13		23		33	
4	SVW38-VPC-065	14		24		34	
5	SVW38-VPC-066DUP	15		25		35	
6	SVW38-VPD-067	16		26		36	
7	SVW38-VPF-068	17		27		37	
8	SVW38-VPG-069	18		28		38	
9	SVW38-VPJ-070	19		29		39	
10	BLK	20		30		40	

Notes:

## Laboratory Data Consultants, Inc. Data Validation Report

**Project/Site Name:** NASA JPL  
**Collection Date:** December 28, 2000  
**LDC Report Date:** March 1, 2001  
**Matrix:** Air  
**Parameters:** Volatile Halogenated/Aromatic Hydrocarbons  
**Validation Level:** EPA Level III  
**Laboratory:** HP Labs

**Sample Delivery Group (SDG):** 2K1228W1

### Sample Identification

SVW37-VPA-071  
SVW37-VPA-072DUP  
SVW37-VPB-073  
SVW37-VPC-074  
SVW37-VPD-075  
SVW37-VPE-076  
SVW37-VPH-077  
SVW37-VPH-078DUP  
SVW37-VPI-079  
SVW37-VPJ-080

## Introduction

This data review covers 10 air samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA SW 846 Methods 8010 and 8020 for Volatile Halogenated/Aromatic Hydrocarbons.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review (October 1999) as there are no current guidelines for the method stated above.

A table summarizing all data qualification flags is provided at the end of this report. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blank results are summarized in Section III.

Field duplicates are summarized in Section IX.

Raw data were not reviewed for this SDG. The review was based on QC data.

The following are definitions of the data qualifiers:

- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- J Indicates an estimated value.
- R Quality control indicates the data is not usable.
- N Presumptive evidence of presence of the constituent.
- UU Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## **I. Technical Holding Times**

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## **II. Calibration**

### **a. Initial Calibration**

Initial calibration of compounds was performed for the primary (quantitation) column and confirmation column as required by these methods.

The percent relative standard deviations (%RSD) were less than or equal to 20.0% for all compounds.

### **b. Calibration Verification**

Calibration verification was performed at the required frequencies.

The percent differences (%D) of calibration factors in continuing standard mixtures were within the 15.0% QC limits.

## **III. Blanks**

Method blanks were reviewed for each matrix as applicable. No volatile halogenated/aromatic hydrocarbon contaminants were found in the method blanks.

## **IV. Accuracy and Precision Data**

### **a. Surrogate Recovery**

Surrogates were added to all samples and blanks as required by the methods. All surrogate recoveries (%R) were within QC limits.

### **b. Matrix Spike/Matrix Spike Duplicates**

Matrix spike (MS) and matrix spike duplicate (MSD) analyses were not required by the method.

### **c. Laboratory Control Samples**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits.

## V. Target Compound Identification

Raw data were not reviewed for this SDG.

## VI. Compound Quantitation and CRQLs

Raw data were not reviewed for this SDG.

## VII. System Performance

Raw data were not reviewed for this SDG.

## VIII. Overall Assessment of Data

Data flags are summarized at the end of this report.

## IX. Field Duplicates

Samples SVW37-VPA-071 and SVW37-VPA-072DUP and samples SVW37-VPH-077 and SVW37-VPH-078DUP were identified as field duplicates. No volatile halogenated/aromatic hydrocarbons were detected in any of the samples with the following exceptions:

Compound	Concentration (ug/L)		RPD
	SVW37-VPH-077	SVW37-VPH-078DUP	
Carbon tetrachloride	3.5	3.2	9
1,1,2-Trichlorotrifluoroethane	1.5	1.2	22

## X. Field Blanks

No field blanks were identified in this SDG.



**NASA JPL**

**Volatile Halogenated/Aromatic Hydrocarbons - Data Qualification Summary - SDG 2K1228W1**

No Sample Data Qualified in this SDG

**NASA JPL**

**Volatile Halogenated/Aromatic Hydrocarbons - Laboratory Blank Data Qualification Summary - SDG 2K1228W1**

No Sample Data Qualified in this SDG

GEOFON PROJECT # 04-4304-480

JPL

4800 OAK GROVE DRIVE

PASADENA, CA

HP Labs Project #2K1228W1

GC SHIMADZU 14A FRONT

VOLATILE HALOGENATED AND AROMATIC HYDROCARBONS (EPA Method 8010/8020) ANALYSES OF SOIL VAPOR

SOIL VAPOR DATA IN UG/L-VAPOR

	BLANK	SVW37-VPA-071	SVW37-VPA-072 DUP	SVW37-VPB-073	SVW37-VPC-074	SVW37-VPD-075
DATE	12/28/00	12/28/00	12/28/00	12/28/00	12/28/00	12/28/00
SAMPLING TIME	06:04	06:52	07:11	07:36	08:01	08:24
ANALYSIS TIME	06:04	06:52	07:16	07:39	08:03	08:27
SAMPLING DEPTH (feet)	--	25	25	40	60	80
VOLUME WITHDRAWN (cc)	200	100	100	160	240	320
VOLUME INJECTED	1	1	1	1	1	1
DILUTION FACTOR	1	1	1	1	1	1
CARBON TETRACHLORIDE	nd	nd	nd	1.4	nd	nd
CHLOROETHANE/BROMOMETHANE	nd	nd	nd	nd	nd	nd
CHLOROFORM	nd	nd	nd	nd	nd	nd
1,1-DICHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,2-DICHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,1-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd
CIS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd
TRANS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd	nd
DICHLOROMETHANE	nd	nd	nd	nd	nd	nd
TETRACHLORO ETHENE	nd	nd	nd	nd	nd	nd
1,1,1,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,1,2,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,1,1-TRICHLORO ETHANE	nd	nd	nd	nd	nd	nd
1,1,2-TRICHLORO ETHANE	nd	nd	nd	nd	nd	nd
TRICHLORO ETHENE	nd	nd	nd	nd	nd	nd
VINYL CHLORIDE	nd	nd	nd	nd	nd	nd
TRICHLOROFLUOROMETHANE (FR11)	nd	nd	nd	nd	nd	nd
DICHLORODIFLUOROMETHANE (FR12)	nd	nd	nd	nd	nd	nd
1,1,2-TRICHLOROTRIFLUOROETHANE (FR113)	nd	nd	nd	nd	nd	nd
BENZENE	nd	nd	nd	nd	nd	nd
ETHYLBENZENE	nd	nd	nd	nd	nd	nd
TOLUENE	nd	nd	nd	nd	nd	nd
m&p-XYLENES	nd	nd	nd	nd	nd	nd
o-XYLENE	nd	nd	nd	nd	nd	nd
SURROGATES						
1,4 DIFLUORO BENZENE	98%	96%	93%	98%	94%	99%
CHLOROBENZENE	92%	95%	93%	97%	94%	99%
4 BROMOFLUORO BENZENE	94%	97%	96%	100%	96%	102%

ND INDICATES NOT DETECTED AT A DETECTION LIMIT OF 1.0 UG/L-VAPOR FOR EACH COMPOUND

ANALYSES PERFORMED ON-SITE IN DOHS CERTIFIED MOBILE LABORATORY (CERT #1745)

ANALYSES PERFORMED BY: ALLEN GLOVER

DATA REVIEWED BY: JAMES E. PICKER

9/3/10



GEOFON PROJECT # 04-4304-480

JPL

4800 OAK GROVE DRIVE

PASADENA, CA

HP Labs Project #2K1228W1

GC SHIMADZU 14A FRONT

VOLATILE HALOGENATED AND AROMATIC HYDROCARBONS (EPA Method 8010/8020) ANALYSES OF SOIL VAPOR

SOIL VAPOR DATA IN UG/L-VAPOR

	SVW37-VPE-076	SVW37-VPH-077	SVW37-VPH-078 DUP	SVW37-VPI-079	SVW37-VPJ-080
DATE	12/28/00	12/28/00	12/28/00	12/28/00	12/28/00
SAMPLING TIME	08:47	09:10	09:35	10:01	10:25
ANALYSIS TIME	08:50	09:15	09:40	10:04	10:28
SAMPLING DEPTH (feet)	100	155	155	170	180
VOLUME WITHDRAWN (cc)	400	620	620	680	720
VOLUME INJECTED	1	1	1	1	1
DILUTION FACTOR	1	1	1	1	1
CARBON TETRACHLORIDE	5.9	3.5	3.2	4.1	3.7
CHLOROETHANE/BROMOMETHANE	nd	nd	nd	nd	nd
CHLOROFORM	nd	nd	nd	nd	nd
1,1-DICHLORO ETHANE	nd	nd	nd	nd	nd
1,2-DICHLORO ETHANE	nd	nd	nd	nd	nd
1,1-DICHLORO ETHENE	nd	nd	nd	nd	nd
CIS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd
TRANS-1,2-DICHLORO ETHENE	nd	nd	nd	nd	nd
DICHLOROMETHANE	nd	nd	nd	nd	nd
TETRACHLORO ETHENE	nd	nd	nd	nd	nd
1,1,1,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd
1,1,2,2-TETRACHLORO ETHANE	nd	nd	nd	nd	nd
1,1,1-TRICHLORO ETHANE	nd	nd	nd	nd	nd
1,1,2-TRICHLORO ETHANE	nd	nd	nd	nd	nd
TRICHLORO ETHENE	1.7	nd	nd	1.4	nd
VINYL CHLORIDE	nd	nd	nd	nd	2.2
TRICHLOROFLUOROMETHANE (FR11)	nd	nd	nd	nd	nd
DICHLORODIFLUOROMETHANE (FR12)	nd	nd	nd	nd	nd
1,1,2-TRICHLOROTRIFLUOROETHANE (FR113)	1.1	1.5	1.2	nd	nd
BENZENE	nd	nd	nd	2.0	3.9
ETHYLBENZENE	nd	nd	nd	nd	nd
TOLUENE	nd	nd	nd	nd	nd
m&p-XYLENES	nd	nd	nd	nd	nd
o-XYLENE	nd	nd	nd	nd	nd
SURROGATES					
1,4 DIFLUORO BENZENE	96%	94%	101%	99%	94%
CHLOROBENZENE	96%	95%	99%	99%	95%
4 BROMOFLUORO BENZENE	99%	96%	102%	101%	99%

ND INDICATES NOT DETECTED AT A DETECTION LIMIT OF 1.0 UG/L-VAPOR FOR EACH COMPOUND

ANALYSES PERFORMED ON-SITE IN DOHS CERTIFIED MOBILE LABORATORY (CERT #1745)

ANALYSES PERFORMED BY: ALLEN GLOVER

DATA REVIEWED BY: JAMES E. PICKER

9/3/201

LDC #: 6038G23 **VALIDATION COMPLETENESS WORKSHEET**  
 SDG #: 2K1228W1   X   EPA Level III        NFESC Level C  
 Laboratory: HP Labs

Date: 2/28/01  
 Page: 1 of 1  
 Reviewer: PT  
 2nd Reviewer:   

**METHOD:** GC Volatile Halogenated/Aromatic Hydrocarbons (EPA SW 846 Method 8010/8020)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: <u>12/28/00</u>
IIa.	Initial calibration	A	% RSD
IIb.	Calibration verification	A	% D
III.	Blanks	A	
IVa.	Surrogate recovery	A	
IVb.	Matrix spike/Matrix spike duplicates	N	
IVc.	Laboratory control samples	FWA	
V.	Target compound identification	N	
VI.	Compound Quantitation and CRQLs	N	
VII.	System Performance	N	
VIII.	Overall assessment of data	N	
IX.	Field duplicates	SW	*P = 1 + 2    D = 7 + 8
X.	Field blanks	N	* = ND

Note: A = Acceptable  
 N = Not provided/applicable  
 SW = See worksheet

ND = No compounds detected  
 R = Rinsate  
 FB = Field blank

D = Duplicate  
 TB = Trip blank  
 EB = Equipment blank

Validated Samples:

AIR

1	SVW37-VPA-071	D	11	BIK	21		31	
2	SVW37-VPA-072DUP	D	12		22		32	
3	SVW37-VPB-073		13		23		33	
4	SVW37-VPC-074		14		24		34	
5	SVW37-VPD-075		15		25		35	
6	SVW37-VPE-076		16		26		36	
7	SVW37-VPH-077	D <sub>1</sub>	17		27		37	
8	SVW37-VPH-078DUP	D <sub>1</sub>	18		28		38	
9	SVW37-VPI-079		19		29		39	
10	SVW37-VPJ-080		20		30		40	

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

LDC #: 6038G23  
SDG #: 2K/228W/

# TARGET COMPOUND WORKSHEET

Page: 1 of 1  
Reviewer: FT  
2nd Reviewer: N

METHOD: VOA (EPA SW 846 Method 8240/8260/8021))

A. Chloromethane*	P. Bromodichloromethane	EE. Ethylbenzene**	TT. 1,2-Dibromoethane	III. n-Butylbenzene
B. Bromomethane	Q. 1,2-Dichloropropane**	FF. Styrene	UU. 1,1,1,2-Tetrachloroethane	JJJ. 1,2-Dichlorobenzene
C. Vinyl chloride**	R. cis-1,3-Dichloropropene	GG. Xylene, total	VV. Isopropylbenzene	KKK. 1,2,4-Trichlorobenzene
D. Chloroethane	S. Trichloroethene	HH. Vinyl acetate	WW. Bromobenzene	LLL. Hexachlorobutadiene
E. Methylene chloride	T. Dibromochloromethane	II. 2-Chloroethylvinyl ether	XX. 1,2,3-Trichloropropane	MMM. Naphthalene
F. Acetone	U. 1,1,2-Trichloroethane	JJ. Dichlorodifluoromethane	YY. n-Propylbenzene	NNN. 1,2,3-Trichlorobenzene
G. Carbon disulfide	V. Benzene	KK. Trichlorofluoromethane	ZZ. 2-Chlorotoluene	OOO. 1,3,5-Trichlorobenzene
H. 1,1-Dichloroethene**	W. trans-1,3-Dichloropropene	LL. Methyl-tert-butyl ether	AAA. 1,3,5-Trimethylbenzene	PPP. trans-1,2-Dichloroethene
I. 1,1-Dichloroethane*	X. Bromoform*	MM. 1,2-Dibromo-3-chloropropane	BBB. 4-Chlorotoluene	QQQ. cis-1,2-Dichloroethene
J. 1,2-Dichloroethene	Y. 4-Methyl-2-pentanone	NN. Diethyl ether	CCC. tert-Butylbenzene	RRR.
K. Chloroform**	Z. 2-Hexanone	OO. 2,2-Dichloropropane	DDD. 1,2,4-Trimethylbenzene	SSS.
L. 1,2-Dichloroethane	AA. Tetrachloroethene	PP. Bromochloromethane	EEE. sec-Butylbenzene	TTT.
M. 2-Butanone	BB. 1,1,2,2-Tetrachloroethane*	QQ. 1,1-Dichloropropene	FFF. 1,3-Dichlorobenzene	UUU.
N. 1,1,1-Trichloroethane	CC. Toluene**	RR. Dibromomethane	GGG. p-Isopropyltoluene	VVV.
O. Carbon tetrachloride	DD. Chlorobenzene*	SS. 1,3-Dichloropropane	HHH. 1,4-Dichlorobenzene	WWW.

\* = System performance check compounds (SPCC) for RF ; \*\* = Calibration check compounds (CCC) for %RSD. ZZZ — 1,1,2-Trichlorotrifluoroethane (FR113)

Notes: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

LDC #: 6038G23  
SDG #: 2K | 228W |

**VALIDATION FINDINGS WORKSHEET**  
**Field Duplicates**

Page: 1 of 1  
Reviewer: FT  
2nd reviewer: N

**METHOD:** GC Volatiles (EPA SW 846 Method 8010/8020)

Y N N/A  
Y N N/A

Were field duplicate pairs identified in this SDG?

Were target compounds detected in the field duplicate pairs?

Compound	Concentration ( <u>ug/L</u> )		RPD
	<u>7</u>	<u>8</u>	
<u>0</u>	<u>3.5</u>	<u>3.2</u>	<u>9</u>
<u>222</u>	<u>1.5</u>	<u>1.2</u>	<u>22</u>

Compound	Concentration ( )		RPD

Compound	Concentration ( )		RPD

Compound	Concentration ( )		RPD